

PTO/SB/08A (10-01)

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				Application Number	10/825,186
				Filing Date	April 16, 2004
				First Named Inventor	Zhang et al.
				Art Unit	1645 1631
				Examiner Name	To Be Assigned Moran
Sheet	1	of	2	Attorney Docket Number	57953/1221 (ZHA01-01)

U.S. PATENT DOCUMENTS					
Examiner Initials ¹	Cite No. ¹	U.S. Patent Document	Publication Date MM-DD-YYYY	Name of Patentee or Applicant of Cited Document	Pages, Columns, Lines, Where Relevant Passages or Relevant Figures Appear
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FOREIGN PATENT DOCUMENTS						
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/M.M./	1	Chen et al., "Fractionation of Peptide with Disulfide Bond for Quantum Mechanical Calculation of Interaction Energy with Molecules," <i>Journal of Chemical Physics</i> 120(2):839-844 (2004)	
/M.M./	2	Chen et al., "Theoretical Method for Full <i>ab initio</i> Calculation of DNA/RNA-Ligand Interaction Energy," <i>Journal of Chemical Physics</i> 120(24):11386-11391 (2004)	
/M.M./	3	Gao et al., "An Efficient Linear Scaling Method for <i>ab initio</i> Calculation of Electron Density of Proteins," <i>Chemical Physics Letters</i> 394:293-297 (2004)	

Examiner Signature	/Marjorie Moran/	Date Considered	03/18/2007
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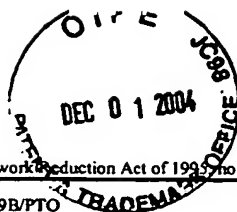
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OTHER PRIOR ART - NON PATENT LITERATURE DOCUMENTS				
Examiner Initials*	Cite No. ¹	Include name of the author (in CAPITAL LETTERS), title of the article (when appropriate), title of the item (book, magazine, journal, serial, symposium, catalog, etc.), date, page(s), volume-issue number(s), publisher, city and/or country where published.		T ²
/M.M./	4	Xiang et al., "Fully Quantum Mechanical Energy Optimization for Protein-Ligand Structure," <i>Journal of Computational Chemistry</i> 25(12):1431-1437 (2004)		
	5	Zhang et al., "Full <i>ab initio</i> Computation of Protein-Water Interaction Energies," <i>Journal of Theoretical and Computational Chemistry</i> 3(1):43-49 (2004)		
	6	Zhang et al., "Molecular Caps for Full Quantum Mechanical Computation of Peptide-Water Interaction Energy," <i>Journal of Computational Chemistry</i> 24(15):1846-1852 (2003)		
	7	Zhang et al., "Molecular Fractionation with Conjugate Caps for Full Quantum Mechanical Calculation of Protein-Molecule Interaction Energy," <i>Journal of Chemical Physics</i> 119(7):3599-3605 (2003)		
	8	Zhang et al., "New Advance in Computational Chemistry: Full Quantum Mechanical <i>ab Initio</i> Computation of Streptavidin-Biotin Interaction Energy," <i>J. Phys. Chem.</i> 107:12039-12041 (2003)		
↓	9	Zhang et al., "Quantum Mechanical Map for Protein-Ligand Binding with Application to β -Trypsin/Benzamidine Complex," <i>Journal of Chemical Physics</i> 120(3):1145-1148 (2004)		

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